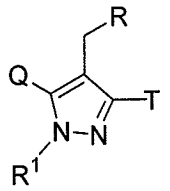
AMENDMENTS TO THE SPECIFICATION

Please replace the paragraphs on pages 17-21 with the following amended paragraphs:

[11] an inhibitor of 1,5-anhydroglucitol/fructose/mannose transporter comprising as an active ingredient a pyrazole derivative represented by the following general formula (I):

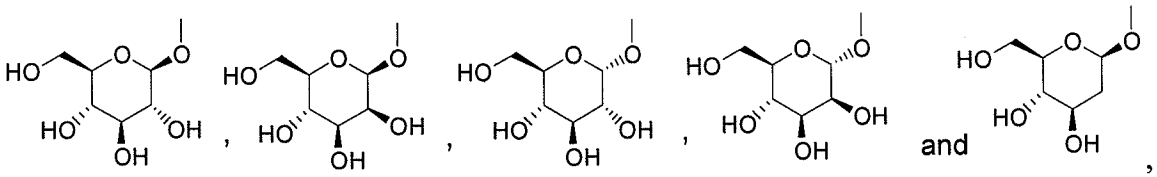


(I)

wherein

R^1 represents a hydrogen atom, a C_{1-6} alkyl group which may have the same or different 1 to 3 groups selected from the following substituent group (A), a C_{2-6} alkenyl group which may have the same or different 1 to 3 groups selected from the following substituent group (A), a C_{2-6} alkynyl group which may have the same or different 1 to 3 groups selected from the following substituent group (A), a C_{3-8} cycloalkyl group which may have the same or different 1 to 3 groups selected from the following substituent group (A), a C_{6-10} aryl group which may have the same or different 1 to 3 groups selected from the following substituent group (B), a C_{2-9} heterocycloalkyl group which may have the same or different 1 to 3 groups selected from the following substituent group (A), or a C_{1-9} heteroaryl group which may have the same or different 1 to 3 groups selected from the following substituent group (B);

one of Q and T represents a group selected from



and the other represents a group represented by the formula: $-(CH_2)_n$-Ar wherein Ar represents a C_{6-10} aryl group which may have the same or different 1 to 3 groups selected from the following substituent group (B) or a C_{1-9} heteroaryl group which may have the same or different 1 to 3 groups selected from the following substituent group (B); and n represents an integral number from 0 to 2, a C_{1-6} alkyl group which may have the same or different 1 to 3 groups selected from the following substituent group (A), a C_{1-6} alkoxy group which may have the same or different 1 to 3 groups selected from the following substituent group (A), an optionally mono or di(C_{1-6} alkyl)-substituted amino group wherein the C_{1-6} alkyl group may have the same or different 1 to 3 groups selected from the following substituent group (A), a C_{3-8} cycloalkyl group which may have the same or different 1 to 3 groups selected from the following substituent group (A), a C_{2-9} heterocycloalkyl group which may have the same or different 1 to 3 groups selected from the following substituent group (A), or a heterocycle-fused phenyl group which may have the same or different 1 to 3 groups selected from the following substituent group (B);

R represents a C_{3-8} cycloalkyl group which may have the same or different 1 to 3 groups selected from the following substituent group (A), a C_{6-10} aryl group which may have the same or different 1 to 3 groups selected from the following substituent group (B), a C_{2-9} heterocycloalkyl group which may have the same or different 1 to 3 groups selected from the following substituent group (A), or a C_{1-9} heteroaryl group which may have the same or different 1 to 3 groups selected from the following substituent group (B);

[substituent group (A)]:

a halogen atom, a nitro group, a cyano group, an oxo group,-G^1, -OG^2, -SG^2, -$N(G^2)_2$, -$C(=O)G^2$, -$C(=O)OG^2$, -$C(=O)N(G^2)_2$, -$S(=O)_2G^2$, -$S(=O)_2OG^2$, -$S(=O)_2N(G^2)_2$, -$S(=O)G^1$, -$OC(=O)G^1$, -$OC(=O)N(G^2)_2$, -$NHC(=O)G^2$, -$OS(=O)_2G^1$, -$NHS(=O)_2G^1$ and –$C(=O)NHS(=O)_2G^1$;

[substituent group (B)]:

a halogen atom, a nitro group, a cyano group, -G^1, -OG^2, -SG^2, -$N(G^2)_2$, -G^3OG^4, -$G^3N(G^4)_2$, -$C(=O)G^2$, -$C(=O)OG^2$, -$C(=O)N(G^2)_2$, -$S(=O)_2G^2$, -$S(=O)_2OG^2$, -$S(=O)_2N(G^2)_2$, -$S(=O)G^1$, -$OC(=O)G^1$, -$OC(=O)N(G^2)_2$, -$NHC(=O)G^2$, -$OS(=O)_2G^1$, -$NHS(=O)_2G^1$ and –$C(=O)NHS(=O)_2G^1$;

in the above substituent group (A) and/or (B),

G^1 represents a C_{1-6} alkyl group which may have the same or different 1 to 3 groups selected from the following substituent group (C), a C_{2-6} alkenyl group which may have the same or different 1 to 3 groups selected from the following substituent group (C), a C_{2-6} alkynyl group which may have the same or different 1 to 3 groups selected from the following substituent group (C), a C_{3-8} cycloalkyl group which may have the same or different 1 to 3 groups selected from the following substituent group (C), a C_{6-10} aryl group which may have the same or different 1 to 3 groups selected from the following substituent group (D), a C_{2-9} heterocycloalkyl group which may have the same or different 1 to 3 groups selected from the following substituent group (C), or a C_{1-9} heteroaryl group which may have the same or different 1 to 3 groups selected from the following substituent group (D);

G^2 represents a hydrogen atom, a C_{1-6} alkyl group which may have the same or different 1 to 3 groups selected from the following substituent group (C), a C_{2-6} alkenyl group which may have the same or different 1 to 3 groups selected from the following substituent group (C), a C_{2-6}

alkynyl group which may have the same or different 1 to 3 groups selected from the following substituent group (C), a C_{3-8} cycloalkyl group which may have the same or different 1 to 3 groups selected from the following substituent group (C), a C_{6-10} aryl group which may have the same or different 1 to 3 groups selected from the following substituent group (D), a C_{2-9} heterocycloalkyl group which may have the same or different 1 to 3 groups selected from the following substituent group (C), or a C_{1-9} heteroaryl group which may have the same or different 1 to 3 groups selected from the following substituent group (D), and with the proviso that G^2 may be the same or different when there are 2 or more G^2 in the substituents;

G^3 represents a C_{1-6} alkyl group;

G^4 represents a C_{1-6} alkyl group which may have the same or different 1 to 3 groups selected from the following substituent group (C), and with the proviso that G^4 may be the same or different when there are 2 or more G^4 in the substituents;

[substituent group (C)]:

a halogen atom, a nitro group, a cyano group, an oxo group, $-G^5$, $-OG^6$, $-SG^6$, $-N(G^6)_2$, $-C(=O)G^6$, $-C(=O)OG^6$, $-C(=O)N(G^6)_2$, $-S(=O)_2G^6$, $-S(=O)_2OG^6$, $-S(=O)_2N(G^6)_2$, $-S(=O)G^5$, $-OC(=O)G^5$, $-OC(=O)N(G^6)_2$, $-NHC(=O)G^6$, $-OS(=O)_2G^5$, $-NHS(=O)_2G^5$ and $-C(=O)NHS(=O)_2G^5$; and

[substituent group (D)]:

a halogen atom, a nitro group, a cyano group, $-G^5$, $-OG^6$, $-SG^6$, $-N(G^6)_2$, $-C(=O)G^6$, $-C(=O)OG^6$, $-C(=O)N(G^6)_2$, $-S(=O)_2G^6$, $-S(=O)_2OG^6$, $-S(=O)_2N(G^6)_2$, $-S(=O)G^5$, $-OC(=O)G^5$, $-OC(=O)N(G^6)_2$, $-NHC(=O)G^6$, $-OS(=O)_2G^5$, $-NHS(=O)_2G^5$ and $-C(=O)NHS(=O)_2G^5$;

in the substituent group (C) and/or (D),

G^5 represents a C_{1-6} alkyl group, a C_{2-6} alkenyl group, a C_{2-6} alkynyl, a C_{3-8} cycloalkyl group, a C_{6-10} aryl group, a C_{2-9} heterocycloalkyl group or a C_{1-9} heteroaryl group; and

G^6 represents a hydrogen atom, a C_{1-6} alkyl group, a C_{2-6} alkenyl group, a C_{2-6} alkynyl, a C_{3-8} cycloalkyl group, a C_{6-10} aryl group, a C_{2-9} heterocycloalkyl group or a C_{1-9} heteroaryl group, and with the proviso that G^6 may be the same or different when there are 2 or more G^6 in the substituents, or a pharmaceutically acceptable salt thereof or a prodrug thereof;